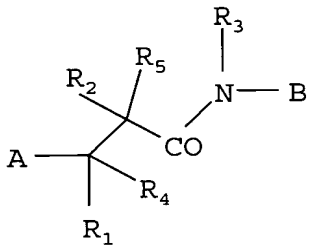
Claims 1-2 (cancelled)

Claim 3 (currently amended): A compound of the formula



(I),

, wherein:

R_1 denotes ~~a hydrogen atom,~~ a C_{1-3}-alkyl ~~or trifluoromethyl~~ group,

R_2 denotes a hydrogen, fluorine, chlorine or bromine atom, a C_{1-3}-alkyl, C_{3-7}-cycloalkyl or C_{1-3}-alkoxy group ~~or, if R_4 and R_5 each denote a hydrogen atom, R_1 and R_2 together denote an n-C_{1-3}-alkylene group optionally substituted by a C_{1-3}-alkyl group~~,

R_3 denotes a hydrogen atom or a C_{1-5}-alkyl group,

R_4 and R_5 ~~each denote a hydrogen atom or~~ together denote another carbon-carbon bond,

A denotes a phenyl~~, naphthyl or tetrahydronaphthyl~~ group substituted by a fluorine, chlorine, bromine or iodine atom, ~~by a C_{1-6}-alkyl, C_{3-7}-cycloalkyl, phenyl, C_{1-3}-alkoxy, cyano, trifluoromethyl or nitro group,~~ whilst the abovementioned monosubstituted phenyl ~~and naphthyl~~ groups ~~may~~ is additionally ~~be~~ substituted by a fluorine, chlorine or bromine atom, ~~by a C_{1-3}-alkyl or C_{1-3}-alkoxy group~~ and the abovementioned disubstituted phenyl groups may additionally be substituted by a C_{1-3}-alkyl or C_{1-3}-alkoxy group, ~~with the proviso that~~

~~A does not denote a phenyl group which is substituted by a halogen atom, by a methyl, pentyl, C_{1-3}-alkoxy or phenyl group or by two C_{1-3}-alkoxy groups, if~~

~~R_3 denotes a hydrogen atom,~~

R_4 and R_5 each denote a hydrogen atom or

R_4 and R_5 together denote another carbon-carbon bond and

B denotes a carboxyphenyl or methoxycarbonylphenyl group,

and A does not denote a phenyl group substituted by a methyl or phenyl group if

R_1 and R_2 each denote a hydrogen atom,

R_3 denotes a hydrogen atom,

R_4 and R_5 together denote another carbon-carbon bond and

B denotes a carboxyphenyl or methoxycarbonylphenyl group,

a naphthyl group,

a chromane or chromene group wherein a methylene group may be replaced by a carbonyl group,

a 5- or 6-membered heteroaryl group optionally substituted in the carbon skeleton by a fluorine, chlorine or bromine atom or by a C_{1-3}-alkyl or C_{1-3}-alkoxy group, whilst the 6-membered heteroaryl groups contain one, two or three nitrogen atoms and the 5-membered heteroaryl groups contain an imino group optionally substituted by a C_{1-3}-alkyl group, an oxygen or sulphur atom or an imino group optionally substituted by a C_{1-3}-alkyl group and an oxygen or sulphur atom or one or two nitrogen atoms and additionally a phenyl ring may be fused to the abovementioned monocyclic heteroaryl groups via two adjacent carbon atoms, whilst said phenyl ring may also be substituted in the carbon skeleton by a fluorine, chlorine or bromine atom, by a C_{1-3}-alkyl or C_{1-3}-alkoxy group,

a phenylvinyl group or

R_1 together with A and the carbon atom between them denote a C_{5-7}-cycloalkylidene group to which a phenyl ring may be fused via two adjacent carbon atoms, whilst said phenyl ring may additionally be substituted by one or two C_{1-3}-alkyl or C_{1-3}-alkoxy groups, whilst the substituents may be identical or different, and

B denotes ~~a 5- or 6-membered heteroaryl group substituted by a carboxy group or by a group which may be converted into a carboxy group *in vivo*,~~

a phenyl ~~or naphthyl~~ group, ~~each of~~ which ~~may be~~is substituted by a carboxy group, ~~by a group which may be converted into a carboxy group *in vivo* or by a group which is negatively charged under physiological conditions,~~ whilst the abovementioned phenyl groups may additionally be substituted

by a fluorine, chlorine, bromine or iodine atom,

by a C_{1-3}-alkyl, trifluoromethyl, ~~phenyl,~~ hydroxy, C_{1-3}-alkoxy, C_{1-3}-alkylsulphonyloxy, phenylsulphonyloxy, carboxy, C_{1-3}-alkoxycarbonyl, formyl, C_{1-3}-alkylcarbonyl, C_{1-3}-alkylsulphonyl, phenylsulphonyl, nitro, pyrrolidino, piperidino, morpholino, N-(C_{1-3}-alkyl)-piperazino, aminoulphonyl, C_{1-3}-alkylaminosulphonyl or di-(C_{1-3}-alkyl)-aminosulphonyl group,

by a C_{1-3}-alkyl group which is substituted by a hydroxy, C_{1-3}-alkoxy, amino, C_{1-4}-alkylamino, di-(C_{1-4}-alkyl)-amino, C_{3-7}-cycloalkylamino, pyrrolidino, piperidino, morpholino, piperazino or N-(C_{1-3}-alkyl)-piperazino group,

by an n-C_{2-3}-alkoxy, C_{2-3}-alkenyl or C_{2-3}-alkynyl group substituted in the 2 or 3 position by a di-(C_{1-3}-alkyl)-amino group,

by an amino group, by an N-(C_{1-3}-alkyl)-amino or N,N-di-(C_{1-3}-alkyl)-amino group wherein the alkyl moiety may in each case be substituted in the 2 or 3 position in relation to the nitrogen atom by a C_{1-3}-alkoxy group, by an N-phenylamino, N-(phenyl-C_{1-3}-alkyl)-amino or N-(pyridyl-C_{1-3}-alkyl)-amino group wherein in each case a hydrogen atom of the abovementioned amino groups may be substituted by a C_{1-3}-alkylsulphonyl, phenyl-C_{1-3}-alkylsulphonyl or phenylsulphonyl group or by a C_{1-7}-alkyl group which may be replaced in the 2 to 5 position by a C_{1-3}-alkoxy, cyano, amino, C_{1-3}-alkylamino, di-(C_{1-3}-alkyl)-amino or tetrazolyl group,

by an aminocarbonyl or C_{1-3}-alkylaminocarbonyl group which may in each case be substituted at the amino-nitrogen atom

by a C_{1-4}-alkyl group which may be substituted by a vinyl, ethynyl, phenyl, pyridyl, imidazolyl, carboxy or trifluoromethyl group or, with the exception of the 2 position relative to the aminocarbonyl nitrogen atom, by a hydroxy, C_{1-3}-alkoxy, C_{1-3}-alkylthio, amino, C_{1-3}-alkylamino, di-(C_{1-3}-alkyl)-amino, C_{1-4}-alkanoylamino or C_{1-5}-alkoxycarbonylamino group,

by a C_{3-7}-cycloalkyl, C_{5-9}-azabicycloalkyl, phenyl, pyridyl, C_{1-3}-alkoxy or di-(C_{1-3}-alkyl)-amino group,

by a C_{1-3}-alkyl group which is substituted by a piperidin-3-yl or piperidin-4-yl group optionally substituted in the 1 position by a C_{1-3}-alkyl or C_{1-5}-alkoxycarbonyl group, or

by an amino, C_{1-3}-alkylamino or phenyl-C_{1-3}-alkylamino group optionally substituted at the amino-nitrogen atom by a C_{1-4}-alkanoyl, C_{1-5}-alkoxycarbonyl, benzoyl, pyrrolidino, piperidino, morpholino or N-(C_{1-3}-alkyl)-piperazino group,

by a carbonyl group substituted by a pyrrolidino, pyrrolino, piperidino, morpholino or N-(C_{1-3}-alkyl)-piperazino group,

by a sulphonyl group substituted by an amino, C_{1-3}-alkylamino, di-(C_{1-3}-alkyl)-amino, pyrrolidino, piperidino, morpholino or N-(C_{1-3}-alkyl)-piperazino group,

by an amino or N-(C_{1-3}-alkyl)-amino group which may in each case be substituted at the amino-nitrogen atom by an aminocarbonyl, C_{1-3}-alkylaminocarbonyl, phenyl-C_{1-3}-alkylaminocarbonyl, phenylaminocarbonyl, phenoxyphenylaminocarbonyl, pyridylaminocarbonyl, pyrrolidinocarbonyl, piperidinocarbonyl, morpholinocarbonyl or N-(C_{1-3}-alkyl)-piperazinocarbonyl group, wherein additionally any hydrogen atom of one of the abovementioned aminocarbonyl groups present may be substituted by a C_{1-3}-alkyl group,

~~by a 5- or 6-membered heteroaryl group,~~

by a dihydro-oxazolyl, dihydro-imidazolyl, 2-oxo-pyrrolidino, 2-oxo-piperidino or 2-oxo-hexamethyleneimino group to which a phenyl ring may be fused via two adjacent carbon atoms,

by an ethynyl group substituted by a phenyl, hydroxymethyl or dimethylamino group,

whilst

additionally the abovementioned mono- or disubstituted phenyl groups may be substituted by another fluorine, chlorine or bromine atom or by one or two other C_{1-3}-alkyl or C_{1-3}-alkoxy groups and two C_{1-3}-alkoxy groups in the o position may be replaced by a methylenedioxy group,

~~and the abovementioned 6-membered heteroaryl groups contain one, two or three nitrogen atoms and the abovementioned 5-membered heteroaryl groups contain an imino group optionally substituted by a C_{1-3}-alkyl group, an oxygen or sulphur atom or an imino group optionally substituted by a C_{1-3}-alkyl group substituted and an oxygen or sulphur atom or one or two nitrogen atoms and additionally a phenyl ring may be fused to the abovementioned monocyclic heteroaryl groups via two adjacent carbon atoms, this phenyl ring optionally being substituted in the carbon skeleton by a fluorine, chlorine or bromine atom or by a C_{1-3}-alkyl or C_{1-3}-alkoxy group, whilst the abovementioned 5-membered monocyclic heteroaryl groups in the carbon skeleton may additionally be substituted by a C_{1-4}-alkyl, trifluoromethyl, phenyl or furanyl group and by another C_{1-3}-alkyl group,~~

and the amino and imino groups mentioned in the definition of the abovementioned groups may additionally be substituted by a group which may be cleaved *in vivo*,

or a physiologically acceptable salt thereof.

Claim 4 (currently amended): A compound of the formula I, according to claim 3, wherein:

B and R_2 to R_5 are defined as in claim 3,

R_1 denotes a ~~hydrogen atom or a~~ C_{1-3}-alkyl group and

A denotes a phenyl~~, naphthyl or tetrahydronaphthyl~~ group substituted by a fluorine, chlorine, bromine or iodine~~ atom or by a C_{1-6}-alkyl, C_{3-7}-cycloalkyl, phenyl, C_{1-3}-alkoxy, trifluoromethyl or nitro group~~, whilst the abovementioned monosubstituted phenyl ~~and naphthyl~~ groups ~~may~~ is additionally ~~be~~ substituted by a fluorine, chlorine or bromine atom~~ or by a C_{1-3}-alkyl or C_{1-3}-alkoxy group, with the proviso that~~

~~A does not denote a phenyl group which may be mono- or disubstituted by halogen atoms, C_{1-4}-alkyl or C_{1-3}-alkoxy groups, wherein the substituents may be identical or different, and does not represent a 4-biphenyl or pentylphenyl group if~~

~~R_1 and R_2 each denote a hydrogen atom or a C_{1-4}-alkyl group,~~
~~R_3 denotes a hydrogen atom,~~
~~R_4 and R_5 each denote a hydrogen atom or~~
~~R_4 and R_5 together denote another carbon-carbon bond and~~
~~B denotes a carboxyphenyl or methoxycarbonylphenyl group,~~

~~a naphthyl group,~~

~~a chromane or chromene group wherein a methylene group may be replaced by a carbonyl group,~~

~~a 5- or 6-membered heteroaryl group optionally substituted in the carbon skeleton by a fluorine, chlorine or bromine atom or by a C_{1-3}-alkyl or C_{1-3}-alkoxy group, whilst the 6-membered heteroaryl groups contain one, two or three nitrogen atoms and the 5-membered heteroaryl groups contain an imino group optionally substituted by a C_{1-3}-alkyl group, an oxygen or sulphur atom or an imino group optionally substituted by a C_{1-3}-alkyl group and an oxygen or sulphur atom or one or two nitrogen atoms and additionally a phenyl ring may be fused to the abovementioned monocyclic heteroaryl groups via two adjacent carbon atoms,~~

~~whilst said phenyl ring may also be substituted in the carbon skeleton by a fluorine, chlorine or bromine atom or by a C_{1-3}-alkyl or C_{1-3}-alkoxy group,~~

or a physiologically acceptable salt thereof.~~the isomers thereof and the salts thereof.~~

Claim 5 (currently amended): A compound of the formula I according to claim 3, wherein:

R_1 denotes ~~a hydrogen atom or~~ a C_{1-3}-alkyl group,

R_2 denotes a hydrogen atom~~or a methyl group or, if R_4 and R_5 each denote a hydrogen atom, R_1 and R_2 together denote a methylene bridge~~,

R_3 denotes a hydrogen atom or a C_{1-5}-alkyl group,

R_4 and R_5 together denote another carbon-carbon bond,

A denotes ~~a phenyl group substituted by a fluorine, chlorine, bromine or iodine atom or by a C_{1-5}-alkyl, cyclohexyl, phenyl, methoxy, cyano or trifluoromethyl group,~~

a phenyl group substituted by fluorine, chlorine or bromine atoms, ~~by methyl or methoxy groups,~~ whilst the substituents may be identical or different, or

a C_{1-3}-alkylphenyl group, which is disubstituted by fluorine, chlorine or bromine atoms, whilst the substituents may be identical or different, ~~with the proviso that~~

~~A does not denote a phenyl group which is substituted by a halogen atom, by a methyl, pentyl, C_{1-3}-alkoxy or phenyl group or by two C_{1-3}-alkoxy groups, if~~

~~R_3 denotes a hydrogen atom,~~
~~R_4 and R_5 each denote a hydrogen atom or~~
~~R_4 and R_5 together denote another carbon-carbon bond and~~
~~B denotes a carboxyphenyl or methoxycarbonylphenyl group,~~

~~and A does not denote a phenyl group which is substituted by a methyl or phenyl group if~~

~~R_1 and R_2 each denote a hydrogen atom,~~

~~R_3 denotes a hydrogen atom,~~

~~R_4 and R_5 together denote another carbon-carbon bond and~~

~~B denotes a carboxyphenyl or methoxycarbonylphenyl group,~~

~~a naphthyl group optionally substituted by a fluorine, chlorine or bromine atom or by a methyl or methoxy group,~~

~~a tetrahydronaphthyl group,~~

~~a chromene group wherein a methylene group is replaced by a carbonyl group,~~

~~a pyridyl, benzofuryl, benzothienyl, quinolyl or isoquinolyl group optionally substituted by a methyl group~~ and

B denotes ~~a cyclohexyl, trimethoxyphenyl, methylenedioxyphenyl, naphthyl, pyridyl, thienyl, pyrazolyl, quinolyl or isoquinolyl group substituted by a carboxy group,~~

a phenyl group substituted by a carboxy, methoxycarbonyl, ethoxycarbonyl, hydroxymethyl, sulpho, tetrazolyl, methylsulphonylaminocarbonyl or phenylsulphonylaminocarbonyl group, which may additionally be substituted

by a fluorine, chlorine, bromine or iodine atom,

by a methyl, trifluoromethyl, ~~phenyl,~~ hydroxymethyl, hydroxy, methoxy, methylsulphonyloxy, 2-dimethylamino-ethoxy, carboxy, nitro, methylsulphonylamino, phenylsulphonylamino, aminosulphonyl, pyrrolidino, piperidino or morpholino group,

by a methyl group which is substituted by an amino, C_{1-3}-alkylamino, cyclopentylamino, pyrrolidino or piperidino group,

by an amino, N-methyl-amino or N-(2-methoxy-ethyl)-amino group which may in each case be substituted at the amino-nitrogen atom

by a C_{1-7}-alkyl or phenyl group,

by an ethyl group which is substituted in the 1 or 2 position by a phenyl or pyridyl group,

by a C_{2-4}-alkyl group which is terminally substituted by a methoxy, cyano, dimethylamino or tetrazolyl group,

by an acetyl, benzoyl, C_{1-5}-alkoxycarbonyl, aminocarbonyl or methylaminocarbonyl group, whilst the aminocarbonyl moiety of the abovementioned groups may in each case additionally be substituted by an optionally phenyl-substituted C_{1-3}-alkyl group, by a phenyl, phenoxyphenyl or pyridyl group,

by a methylsulphonyl, phenylsulphonyl or benzylsulphonyl group,

by an aminocarbonyl or methylaminocarbonyl group which may in each case be substituted at the amino-nitrogen atom

by a C_{1-4}-alkyl, C_{3-6}-cycloalkyl, phenyl, benzyl, pyridyl, pyridylmethyl or methoxy group,

by a methyl group which is substituted by a vinyl, ethynyl, trifluoromethyl, C_{7-9}-azabicycloalkyl, carboxy or imidazolyl group or by a piperidin-4-yl group optionally substituted in the 1 position by a methyl or C_{1-5}-alkoxycarbonyl group,

by a straight-chain or branched C_{2-3}-alkyl group substituted in the 2 or 3 position by a hydroxy, methoxy, methylthio, amino, acetylamino, C_{1-5}-alkoxycarbonylamino, carboxy-, C_{1-5}-alkoxycarbonyl or dimethylamino group,

by a pyrrolidino, piperidino, morpholino, 4-methyl-piperazino, amino or methylamino group, whilst the abovementioned amino and methylamino groups may each additionally be substituted at the amino-nitrogen atom by a methyl, acetyl, benzoyl or C_{1-5}-alkoxycarbonyl group,

by a dihydro-oxazolyl, dihydro-imidazolyl, 2-oxo-pyrrolidino, 2-oxo-piperidino or 2-oxo-hexamethyleneimino group to which a phenyl ring may be fused via two adjacent carbon atoms,

~~by an imidazolyl or 4-methyl-imidazolyl group optionally substituted by a methyl, ethyl or phenyl group, to which a phenyl ring may additionally be fused via two adjacent carbon atoms,~~

~~a pyrazolyl group optionally substituted by a C_{1-4}-alkyl or furanyl group, which may additionally be substituted by a methyl or trifluoromethyl group,~~

by an ethynyl group substituted by a phenyl, hydroxymethyl or dimethylamino group, whilst

additionally the abovementioned mono- or disubstituted phenyl groups may be substituted by another fluorine, chlorine or bromine atom or by one or two other methyl or methoxy groups,

or a physiologically acceptable salt thereof.

Claim 6 (currently amended): A compound of the formula I according to claim 3, wherein:

R_1 denotes ~~a hydrogen atom or~~ a C_{1-3}-alkyl group,

R_2 denotes a hydrogen ~~atom or R_1 and R_2 together denote a methylene group, if R_4 and R_5 each simultaneously denote a hydrogen atom~~,

R_3 denotes a hydrogen atom,

R_4 and R_5 together denote another carbon-carbon bond,

A denotes a phenyl ~~or naphthyl~~ group mono- or disubstituted by a fluorine, chlorine, bromine or iodine atom ~~or by a C_{1-6}-alkyl, C_{3-7}-cycloalkyl or trifluoromethyl group~~, whilst the substituents may be identical or different, ~~with the proviso that~~

~~A does not denote a phenyl group which may be mono- or di-substituted by halogen atoms or C_{1-4}-alkyl groups, wherein the substituents may be identical or different, and does not denote a 4-biphenyl or pentylphenyl group if~~

~~R_1 denotes a hydrogen atom or a C_{1-3}-alkyl group,~~
~~R_2 denotes a hydrogen atom,~~
~~R_3 denotes a hydrogen atom,~~
~~R_4 and R_5 each denote a hydrogen atom or~~
~~R_4 and R_5 together denote another carbon-carbon bond and~~
~~B denotes a carboxyphenyl or methoxycarbonylphenyl group,~~

~~a naphthyl group,~~

~~a chromene group wherein a methylene group is replaced by a carbonyl group,~~

~~a benzothienyl group~~ and

B denotes a phenyl~~, naphthyl, thienyl or pyridinyl group, each of~~ which is substituted by a carboxy group, whilst the abovementioned phenyl groups may additionally be substituted

by a fluorine, chlorine or bromine atom,

by a C_{1-3}-alkyl, hydroxy, C_{1-3}-alkoxy, C_{1-3}-alkylsulphonyloxy, pyrrolidino, piperidino, morpholino or N-(C_{1-3}-alkyl)-piperazino group,

by an n-C_{2-3}-alkoxy, C_{2-3}-alkenyl or C_{2-3}-alkynyl group substituted in the 2 or 3 position by a di-(C_{1-3}-alkyl)-amino group,

by an N-methyl-N-(n-C_{2-3}-alkyl)-amino group substituted in the 2 or 3 position by a di-(C_{1-3}-alkyl)-amino group,

by a di-(C_{1-3}-alkyl)-amino group,

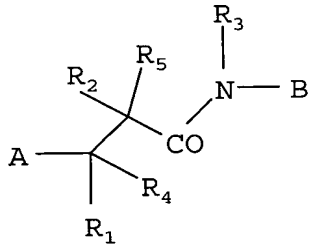
~~by an imidazolyl or pyrazolyl group optionally substituted by a C_{1-4}-alkyl group,~~

by a C_{1-4}-alkylaminocarbonyl, N-(pyridinylmethyl)-aminocarbonyl, pyrrolidinoaminocarbonyl or piperidinoaminocarbonyl group and

may additionally be substituted by another fluorine atom, by another C_{1-3}-alkyl or C_{1-3}-alkoxy group,

or a physiologically acceptable salt thereof.

Claim 7 (currently amended): A compound of the formula I



(I),

~~according to claim 3,~~ wherein:

R_1 denotes a methyl group,

R_2 denotes a hydrogen atom,

R_3 denotes a hydrogen atom,

R_4 and R_5 together denote another carbon-carbon bond,

A denotes a phenyl group substituted by two chlorine or bromine atoms or by a chlorine atom and a bromine atom, ~~a naphthyl, 2-oxo-chromene or benzothienyl group, with the proviso that~~

~~A does not denote a phenyl group disubstituted by halogen atoms if~~

~~R_1 denotes a methyl group,~~
~~R_2 denotes a hydrogen atom,~~
~~R_3 denotes a hydrogen atom,~~
~~R_4 and R_5 each denote a hydrogen atom or~~
~~R_4 and R_5 together denote another carbon-carbon bond and~~
~~B denotes a carboxyphenyl or methoxycarbonylphenyl group,~~

and B denotes a 2-carboxy-phenyl, ~~2-carboxy-thienyl or 2-carboxy-pyridinyl group,~~ whilst the abovementioned 2-carboxy-phenyl group may additionally be substituted in the phenyl nucleus

by a fluorine, chlorine or bromine atom,

by a C_{1-3}-alkyl, hydroxy, C_{1-3}-alkoxy, C_{1-3}-alkylsulphonyloxy or morpholino group,

by an n-C_{2-3}-alkoxy group substituted in the 2 or 3 position by a di-(C_{1-3}-alkyl)-amino group,

by an N-methyl-N-(n-C_{2-3}-alkyl)-amino group substituted in the 2 or 3 position by a di-(C_{1-3}-alkyl)-amino group,

~~by an imidazolyl or pyrazolyl group optionally substituted by a C_{1-4}-alkyl group,~~

by a C_{1-4}-alkylaminocarbonyl, N-(pyridinylmethyl)-aminocarbonyl, pyrrolidinoaminocarbonyl or piperidinoaminocarbonyl group and

may additionally be substituted by another fluorine atom or by another methoxy group,

or a physiologically acceptable salt thereof.

Claims 8 (cancelled)

Claim 9 (original): A pharmaceutical composition containing a compound according to claim 3 together with one or more inert carriers and/or diluents.

Claim 10 (new): A compound selected from the group consisting of:

(a) trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-5-chloro-phenyl)-amide,
(b) trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-phenyl)-amide,
(c) trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-4-methyl-phenyl)-amide,
(d) trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-4,5-dimethoxy-phenyl)-amide,
(e) trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-4-methoxy-5-methyl-phenyl)-amide,
(f) trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(4-bromo-2-carboxy-6-methyl-phenyl)-amide,
(g) trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-5-nitro-phenyl)-amide,
(h) trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-4-iodo-phenyl)-amide,
(i) trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-4-chlorophenyl)-amide,
(j) trans-3-(3,4-dichlorophenyl)-pent-2-enoic acid-N-(2-carboxy-4,5-dimethoxy-phenyl)-amide,
(k) trans-3-(3,4-dichloro-phenyl)-but-2-enoic acid-N-[2-carboxy-4-(morpholin-4-yl)-phenyl]-amide,
(l) trans-3-(3,4-dichloro-phenyl)-but-2-enoic acid-N-(2-carboxy-4-dimethylamino-phenyl)-amide,
(m) trans-3-(2,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-4-fluoro-phenyl)-amide,
(n) trans-3-(2,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-4,5-dimethoxy-phenyl)-amide,
(o) trans-3-(3,4-dichloro-phenyl)-but-2-enoic acid-N-(2-carboxy-4-fluoro-phenyl)-amide,
(p) trans-3-(3,4-difluorophenyl)-but-2-enoic acid-N-(2-carboxy-phenyl)-amide,
(q) trans-3-(3,4-dibromophenyl)-but-2-enoic acid-N-(2-carboxy-phenyl)-amide,
(r) trans-3-(3,4-dibromophenyl)-but-2-enoic acid-N-(2-carboxy-4,5-dimethoxy-phenyl)-amide,
(s) trans-3-(3,4-dibromophenyl)-but-2-enoic acid-N-(2-carboxy-4-methoxy-5-methyl-phenyl)-amide,
(t) trans-3-(3,5-dibromo-4-ethylphenyl)-but-2-enoic acid-N-(2-carboxy-phenyl)-amid,
(u) trans-3-(3-bromo-4-chlorophenyl)-but-2-enoic acid-N-(2-carboxy-phenyl)-amide,

(v) trans-3-(3-chloro-4-bromophenyl)-but-2-enoic acid-N-(2-carboxy-phenyl)-amide,

(w) trans-3-(3-bromo-4-chlorophenyl)-but-2-enoic acid-N-(2-carboxy-4,5-dimethoxy-phenyl)-amide, and

(x) trans-3-(3-chloro-4-bromophenyl)-but-2-enoic acid-N-(2-carboxy-4,5-dimethoxy-phenyl)-amide,

or a physiologically acceptable salt thereof.

Claim 11 (new): Trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-4,5-dimethoxy-phenyl)-amide or a physiologically acceptable salt thereof.

Claim 12 (new): Trans-3-(3,4-dichlorophenyl)-but-2-enoic acid-N-(2-carboxy-4-methoxy-5-methyl-phenyl)-amide or a physiologically acceptable salt thereof.